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## Erratum: Determination of Structure Parameters in Strong-Field Tunneling Ionization Theory of Molecules (Physical Review a (2010) 81 (033423))

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**Erratum: Determination of structure parameters in strong-field tunneling ionization theory of molecules [Phys. Rev. A **81**, 033423 (2010)]**

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There are several errors in Tables V and VI of our article. In Table V, the  $C_{2m}$  of the HOMO-1 ( $1\pi$ ) of CO molecule should be 0.014. In Table VI, the binding energies of  $2p\pi_g$  and  $2p\pi_u$  of  $\text{H}_2^+$  should be 0.2267 and 0.4288, respectively.